

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071216\
 Data File : BG023053.D
 Acq On : 13 Jul 2016 2:47
 Operator : UM/SJ
 Sample : H3980-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FB-070716

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG070816.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.217	400	405	411	rBV	131298	213450	1.86%	0.434%
2	5.693	480	486	497	rBV	1122336	1858618	16.22%	3.777%
3	7.303	753	760	768	rBV	869170	1511931	13.19%	3.072%
4	7.679	816	824	834	rBV	1971863	3438074	30.00%	6.986%
5	8.149	897	904	912	rVB	360307	634921	5.54%	1.290%
6	8.478	952	960	968	rBV	1581364	2783501	24.29%	5.656%
7	9.324	1094	1104	1113	rBV	1611017	2951006	25.75%	5.997%
8	10.981	1378	1386	1398	rVB	581564	1068790	9.33%	2.172%
9	13.420	1793	1801	1809	rBV	4815600	7512367	65.55%	15.265%
10	14.800	2028	2036	2046	rBV2	1127133	1849017	16.13%	3.757%
11	16.293	2282	2290	2304	rBV	4328804	6723704	58.67%	13.663%
12	17.556	2498	2505	2512	rBV2	1445815	2222759	19.40%	4.517%
13	20.159	2941	2948	2956	rBV	8035513	11460109	100.00%	23.287%
14	21.863	3230	3238	3245	rBV	1488437	2598309	22.67%	5.280%
15	25.253	3803	3815	3827	rBV	750813	2385305	20.81%	4.847%

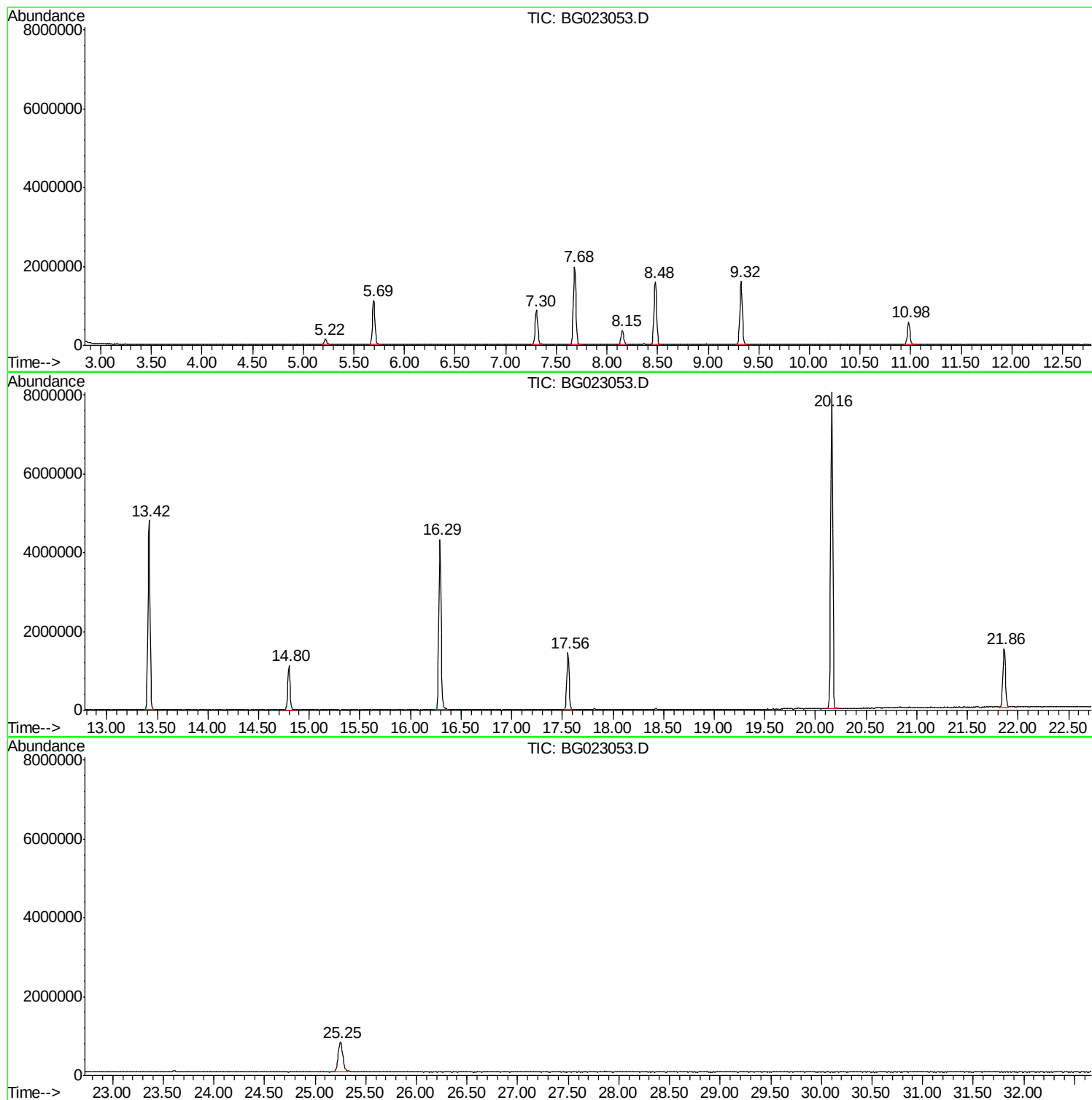
Sum of corrected areas: 49211861

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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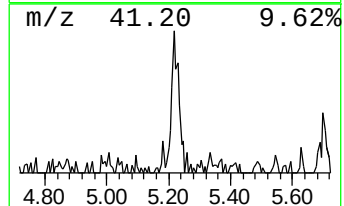
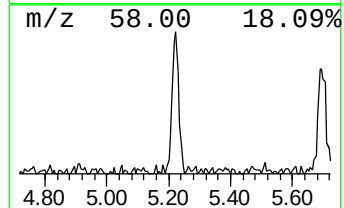
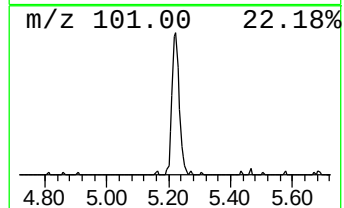
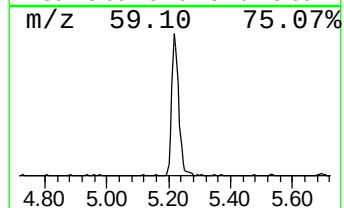
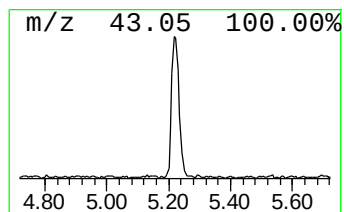
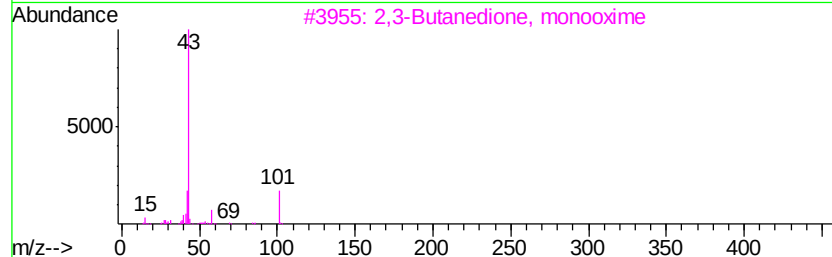
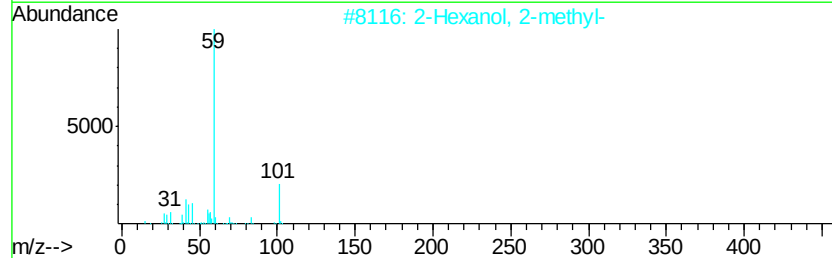
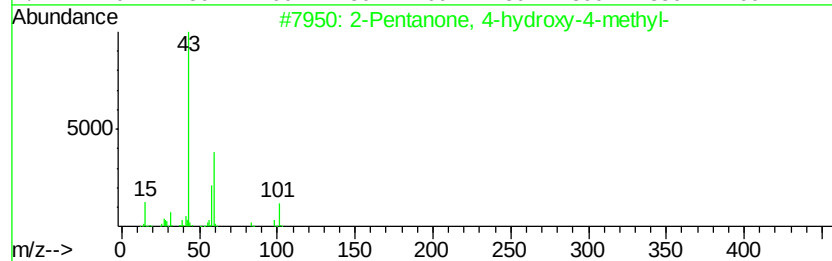
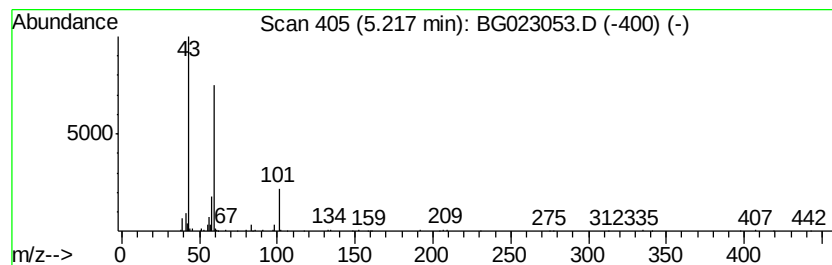
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	6.72 ng	213450	1,4-Dichlorobenzene-d4	8.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	36
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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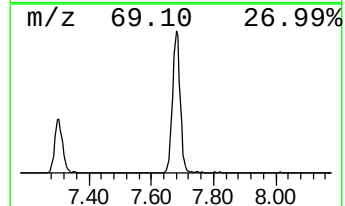
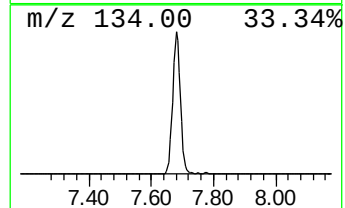
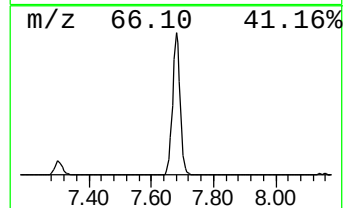
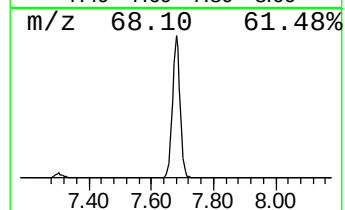
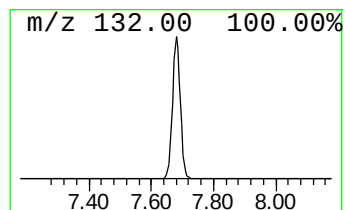
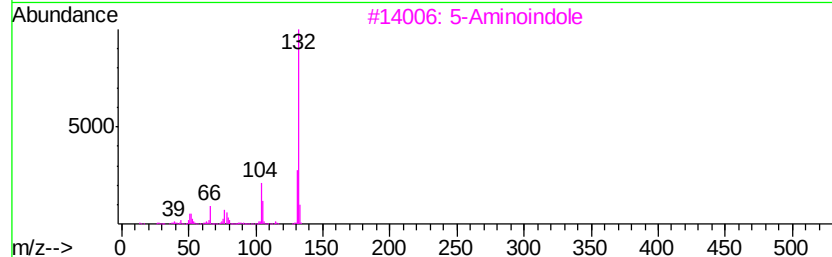
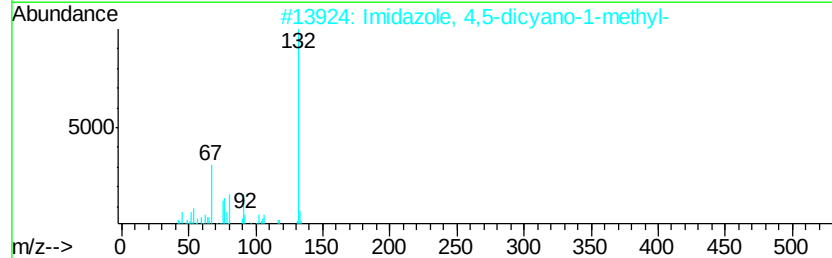
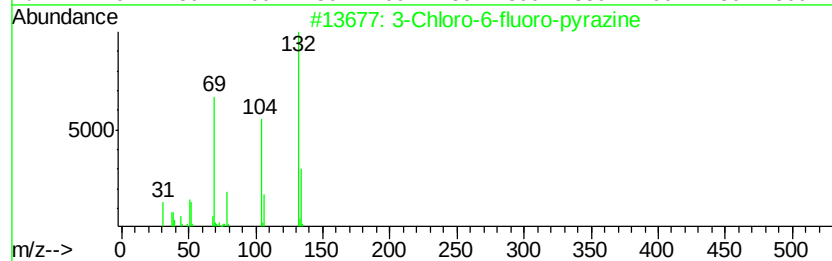
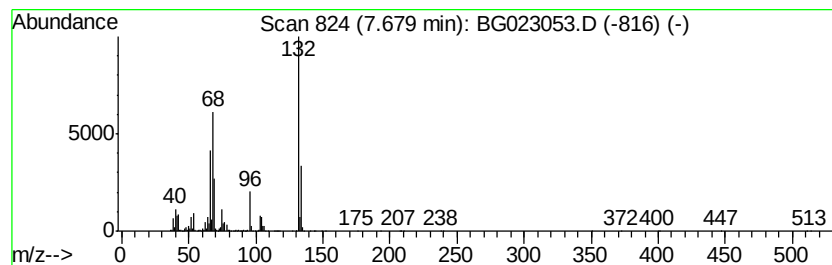
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	108.30 ng	3438070	1,4-Dichlorobenzene-d4	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
2		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	11
3		5-Aminoindole	132	C8H8N2	005192-03-0	9
4		1,2,3-Trifluorobenzene	132	C6H3F3	001489-53-8	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	6.7	ng	213450	1	8.15	634921	20.0
unknown7.68	7.68	108.3	ng	3438070	1	8.15	634921	20.0